

EXHIBIT B

Presentation by Dr. James C. Stevens of The Dow Chemical Company

The Use of High Throughput Experimentation In Catalyst Research

J. C. Stevens, R. Rosen

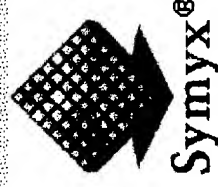
The Dow Chemical Company
Freeport, TX

T. Boussie, G. Diamond, C. Goh, K. Hall, A. M. LaPointe, M. Leclerc,
C. Lund, V. Murphy, J. Shoemaker, H. Turner, U. Tracht, T. Uno

Symyx Technologies
Santa Clara, CA

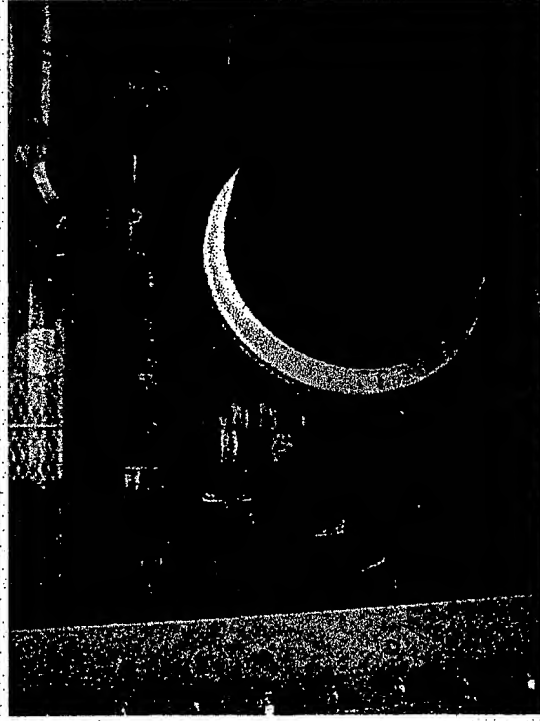


FLEXPO 2002

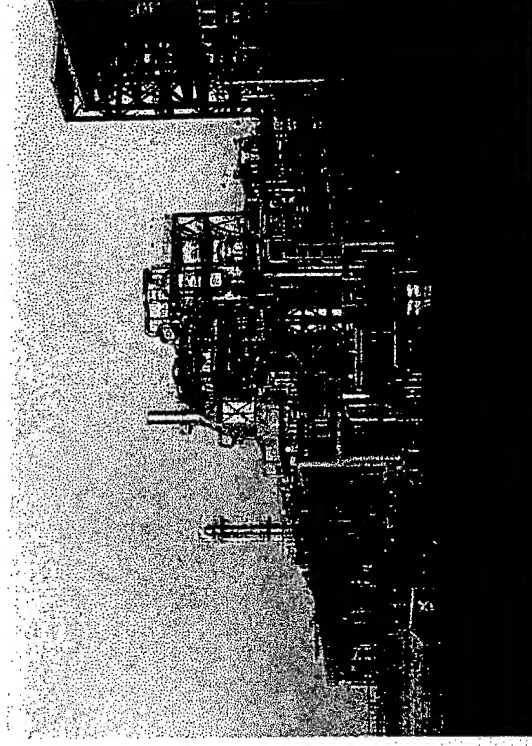


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New Polyolefin Catalysts Drive New Business Growth

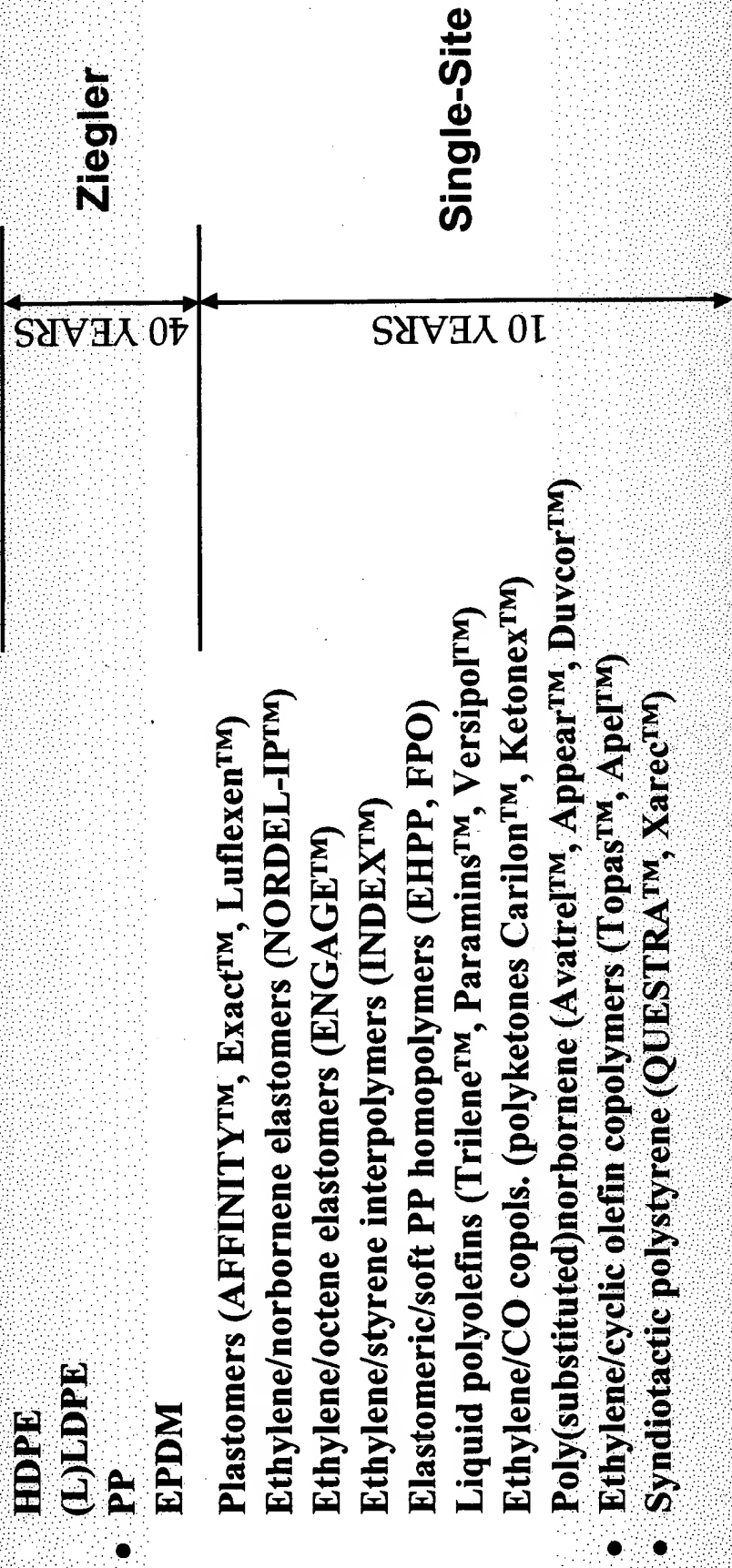


From this



To this

The Rate of New Product Creation Has Been Faster With Single-Site Catalysts



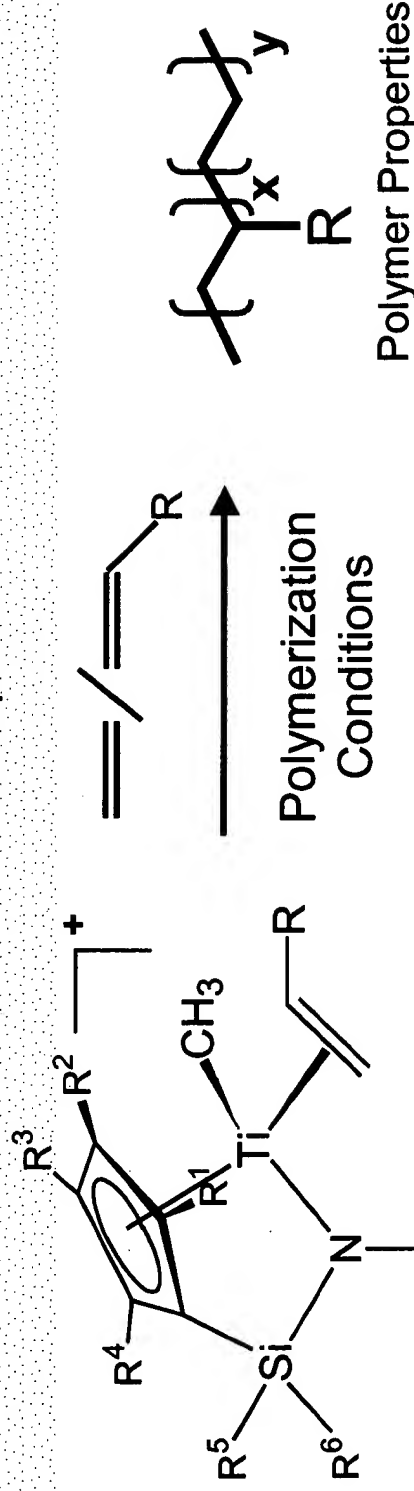
Rate of New Market Creation. Source: STA* research

Single-Site Catalyst Discovery Research Centers On :

- ❖ **More Productive Catalysts**
- ❖ **Higher Temperature Operation**
- ❖ **Higher Molecular Weights**
- ❖ **Higher Comonomer Contents**
- ❖ **Improved / Novel Control of Sequence Distributions**
- ❖ **Incorporation of New Comonomers**
- ❖ **Improved / Differentiated Products**
- ❖ **Novel Polymer Microstructures and Tacticity**

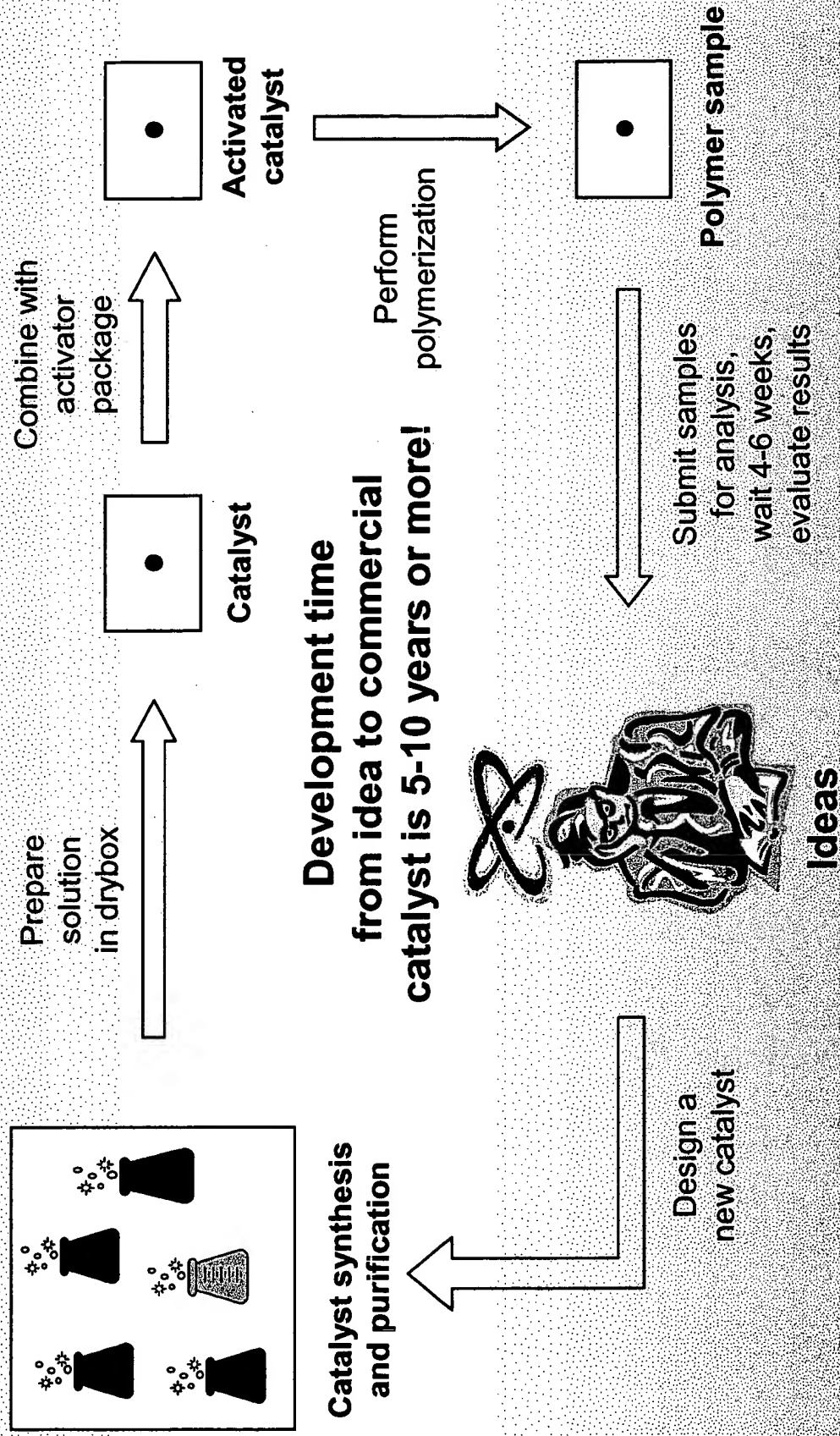
Complete Evaluation of the Structure / Activity Relationships for any Catalyst Family is a Daunting Task

The Energy Surface for a Copolymerization is Highly Complex

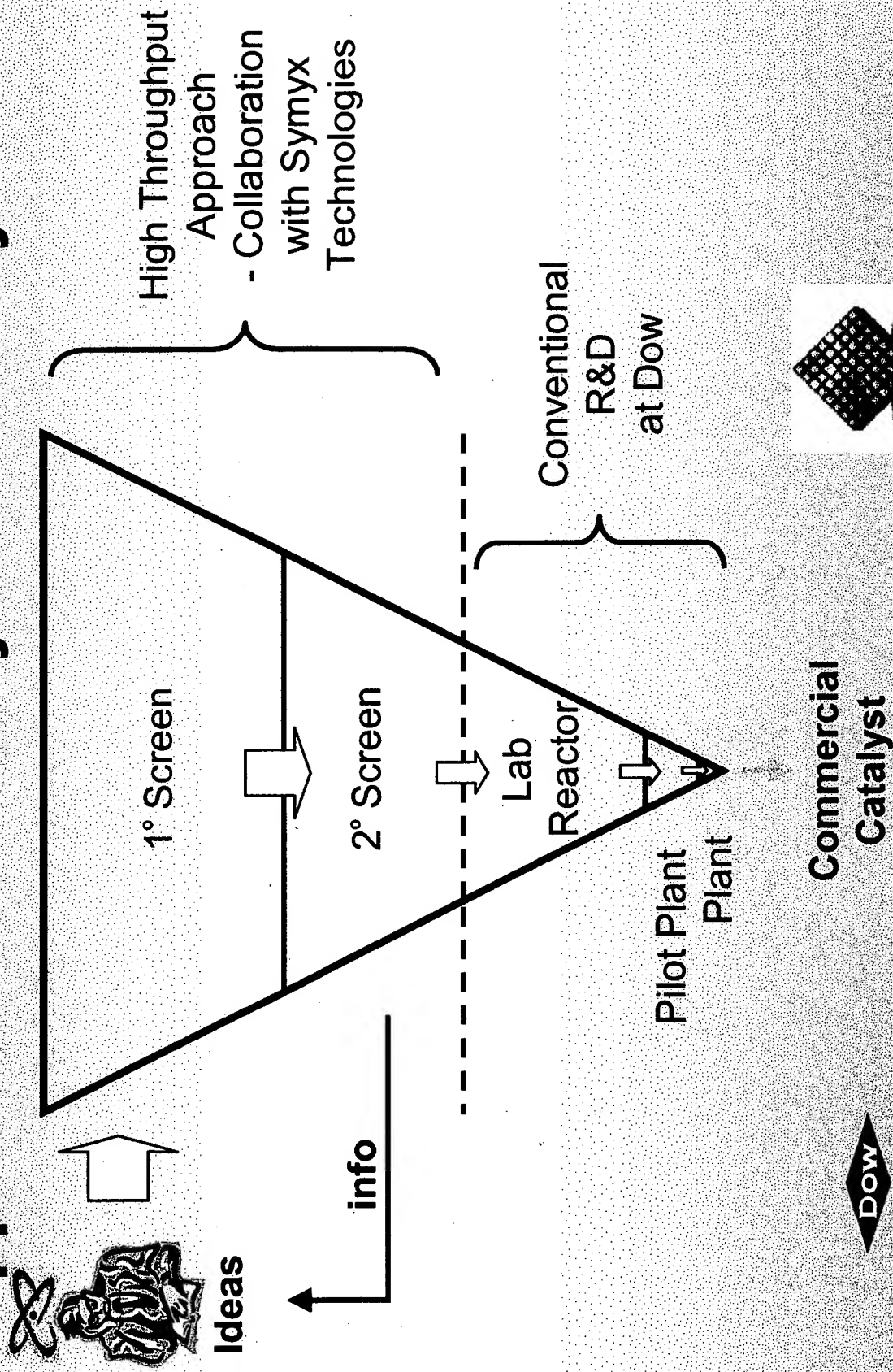


- 1) Properties of active site and polymer can be affected *dramatically and unpredictably* by variations in choice of Metal, Ligand(s), Activation Method and Process Conditions
- 2) Choosing Ten Substituents for Each Position on this ligand Generates around 10,000,000 Compounds !!!

The Conventional Way of Doing Catalyst R&D is Effective But Can Be Slow

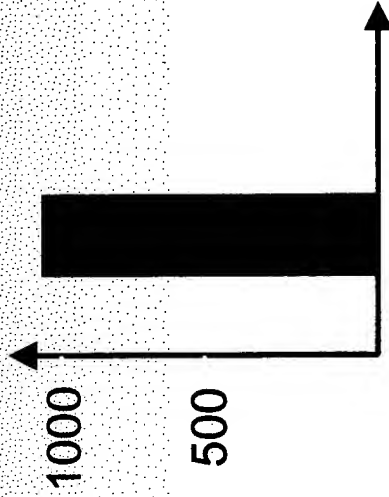


Combinatorial and High-Throughput Approaches to Catalyst Discovery



Experimental Throughput of Symyx / Dow Collaboration

Experiments per Day



**Primary
Screening**

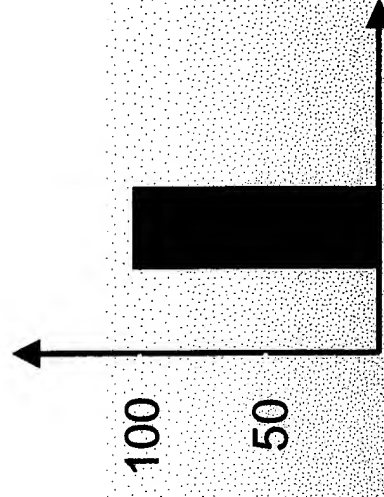
“Discovery Screening”

High Throughput

Lower Precision

Yes/No Approach

Hits and Information



**Secondary
Screening**

“Optimization Screening”

Moderate Throughput

Higher Precision

“Experiment” means catalyst synthesis,
polymerization, and analysis of performance,
polymer properties, etc.

High Throughput Catalyst Discovery Requires:

Large Number of Ligands with Suitable Diversity

Efficient Methods of Attaching Ligands to Metals

Comprehensive Activation Strategies

High Throughput **Primary** Screen

- Rapid analytical for activity, molecular weight, at smaller scales

High Throughput **Secondary** Screen

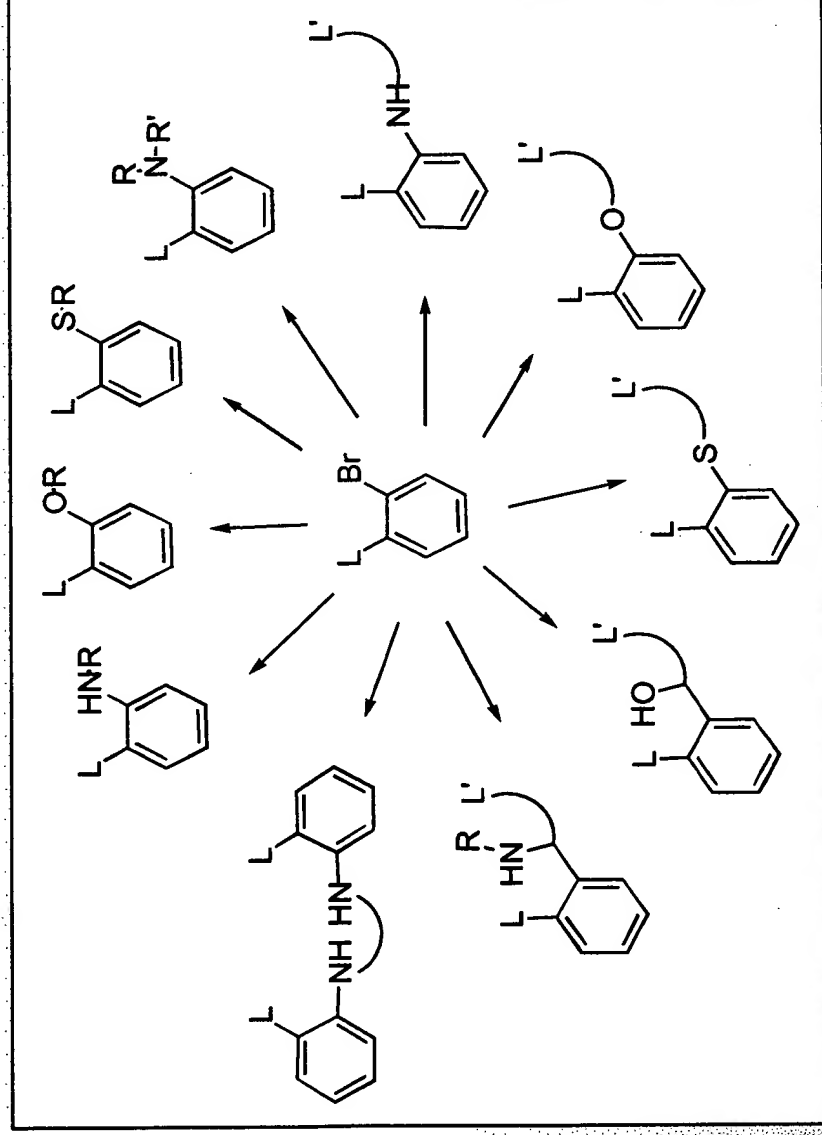
- Rapid analytical for activity, molecular weight, MWD, comonomer incorporation, tacticity, etc

Ligand Diversity: The Building Block Strategy

How can we exploit parallel synthesis techniques to create *diverse* ligand libraries?

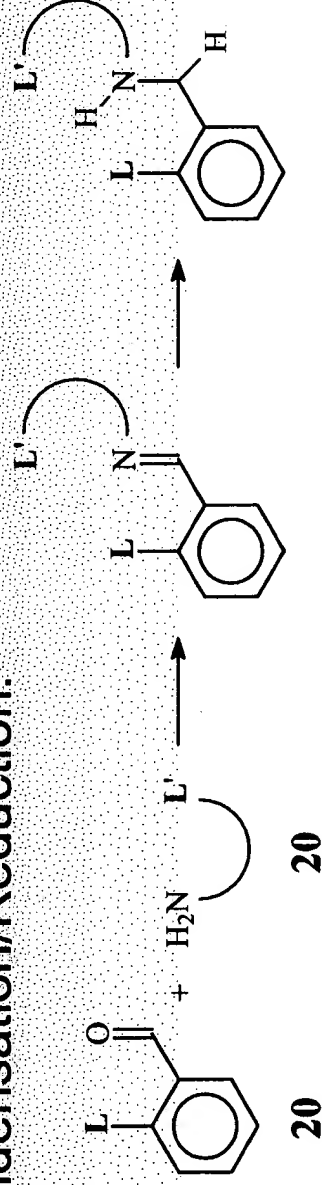
- Focus on post-metallocene catalyst systems
- Create diverse functionalized building blocks
- Connect building blocks using high-yield coupling reactions

The Parallel Building-Block Approach Leads To Significant Diversity:

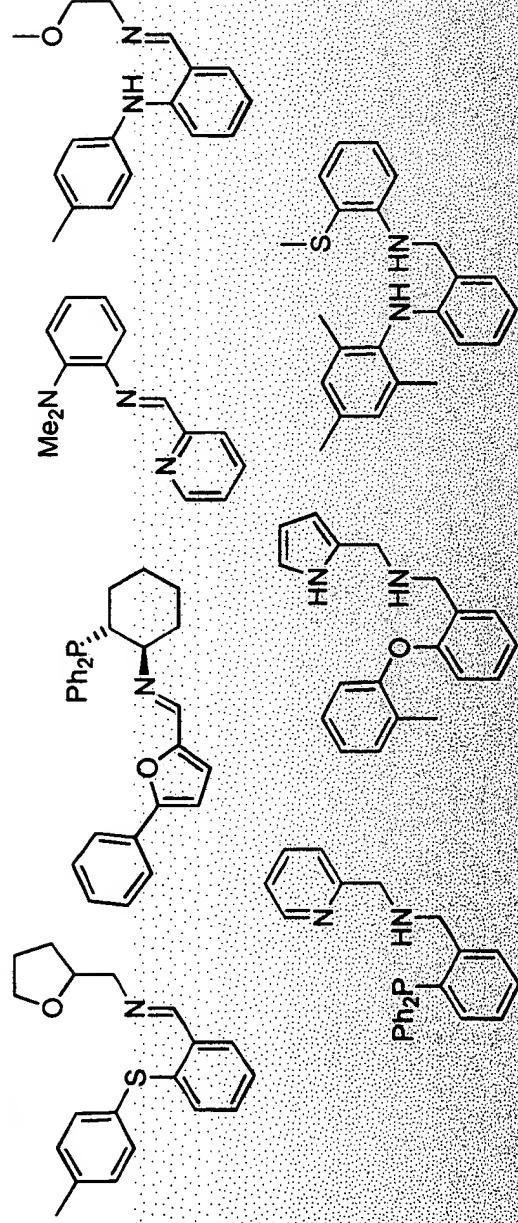


One Example of Ligand Diversity via General Transformations

imine Condensation/Reduction:



20 x 20 x 2 = 800 Examples !!



800 ligands x 6 metals x 3
monomers x 5 activators x
3 temperatures = 216,000
experiments!

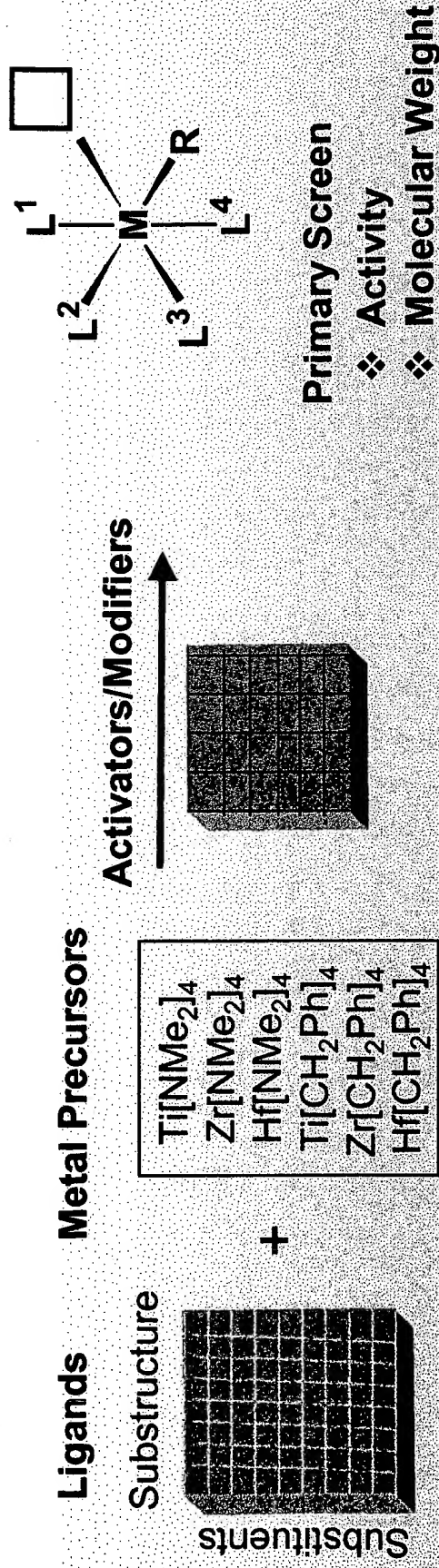
Issues With Parallel Synthesis of Activated Metal Complexes

How can I rapidly ...

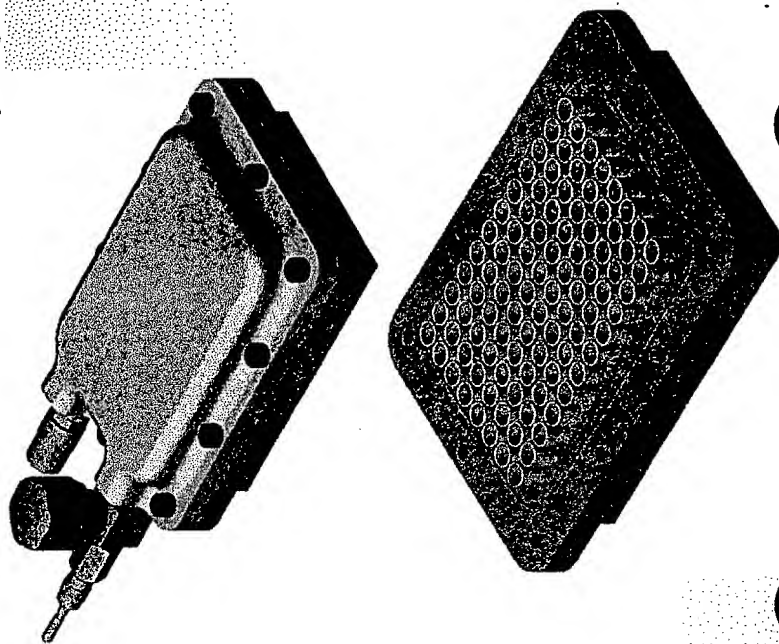
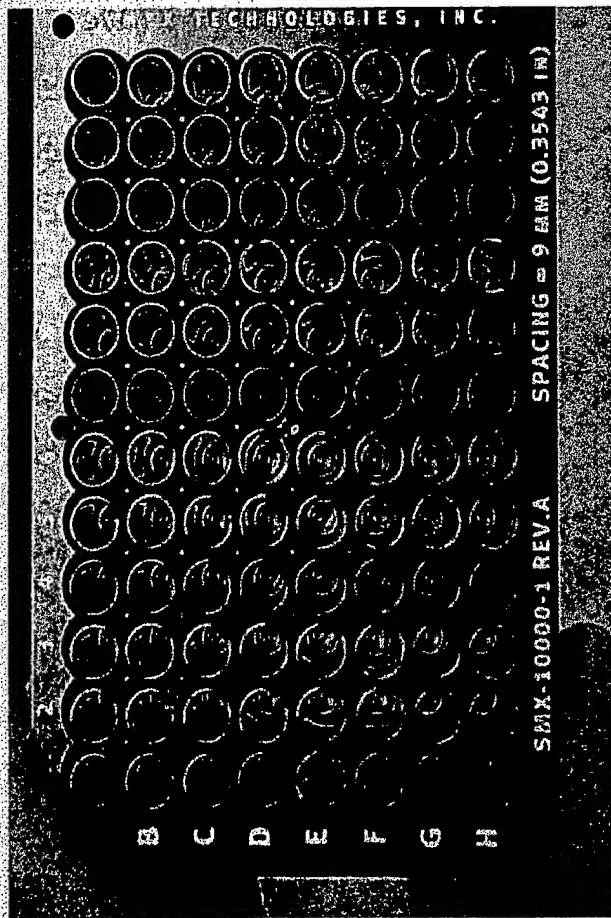
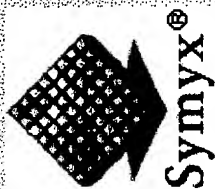
- Synthesize a diverse set of organometallic compounds, and
- know what I made, and
- select the appropriate activator?

By ...

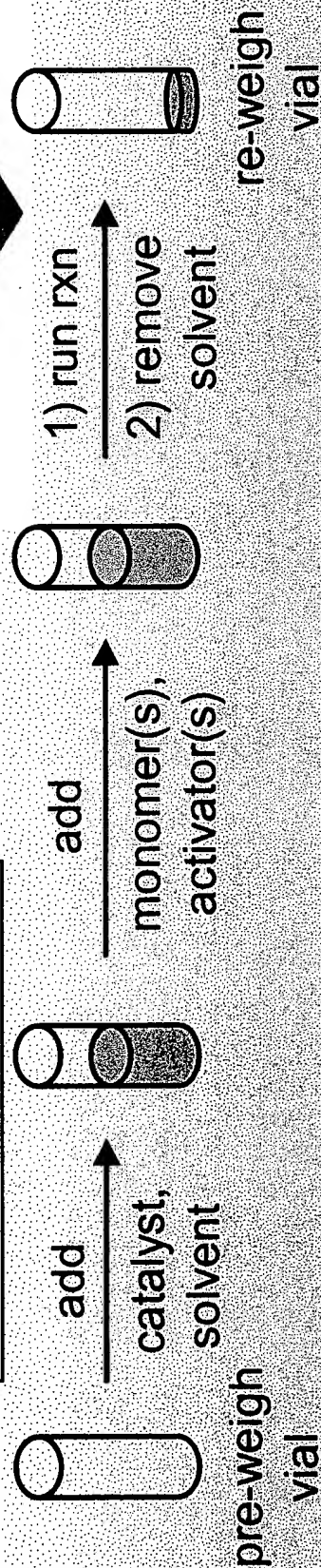
- accommodating multiple organometallic routes, and
- multiple activation chemistries for each metal / ligand set, and
- validating chemistry with models before experiment and for all hits



Rapid Primary Screen Using Post Reaction Gravimetric Analysis

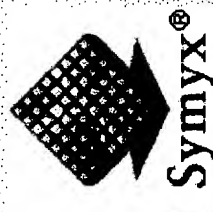
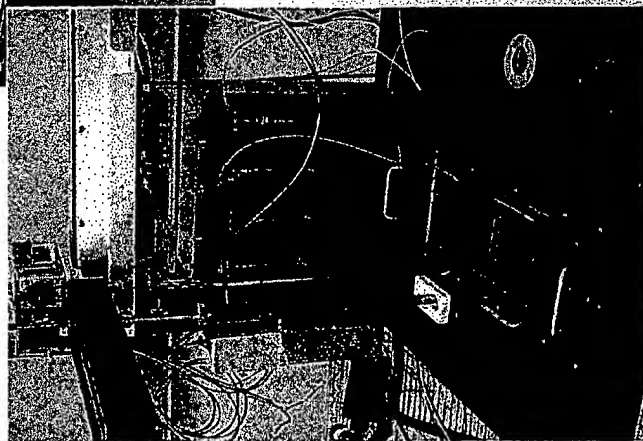
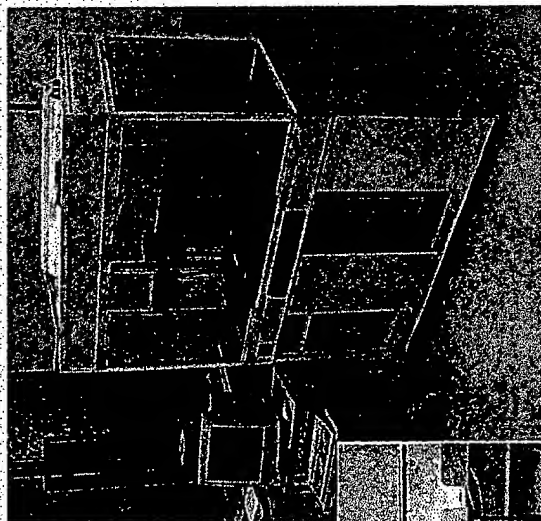


All Procedures Employ
Robotic Liquid Handling

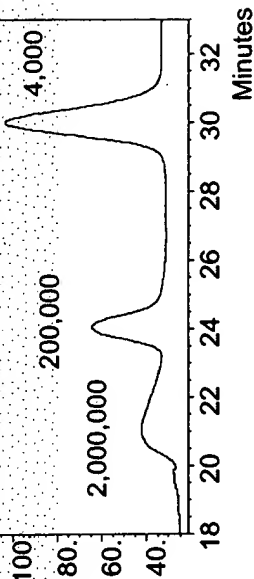


High-Throughput Molecular Weight Measurements

Symyx Rapid GPC™

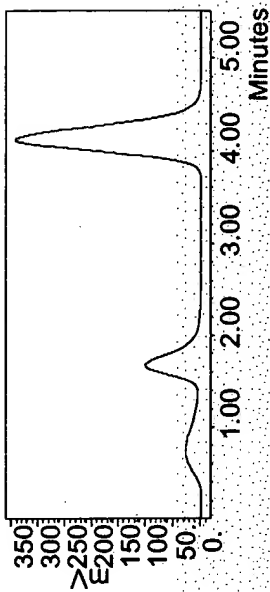


Conventional (PL-GPC210)



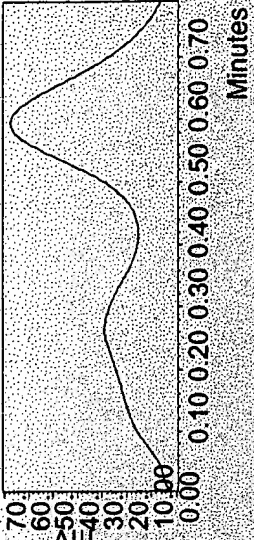
**Conventional:
40 min/sample**

High Resolution



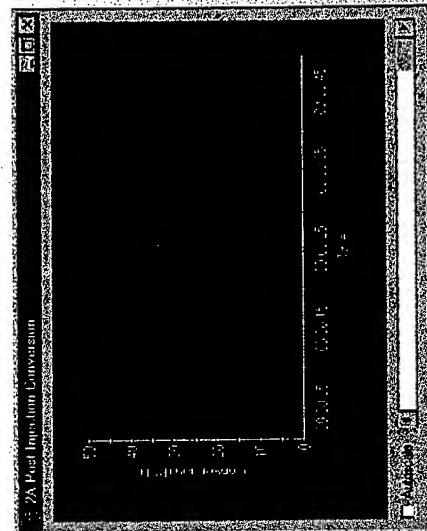
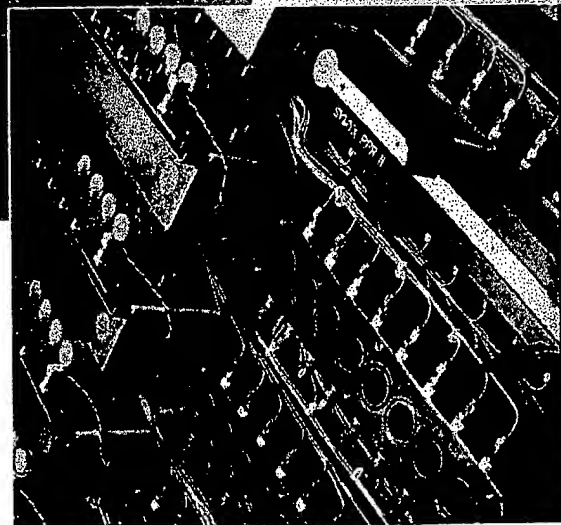
6 Minutes

High Throughput



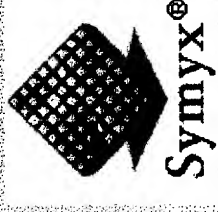
90 Seconds

Secondary Screening Tools Symyx Parallel Polymerization Reactor (PPR®)



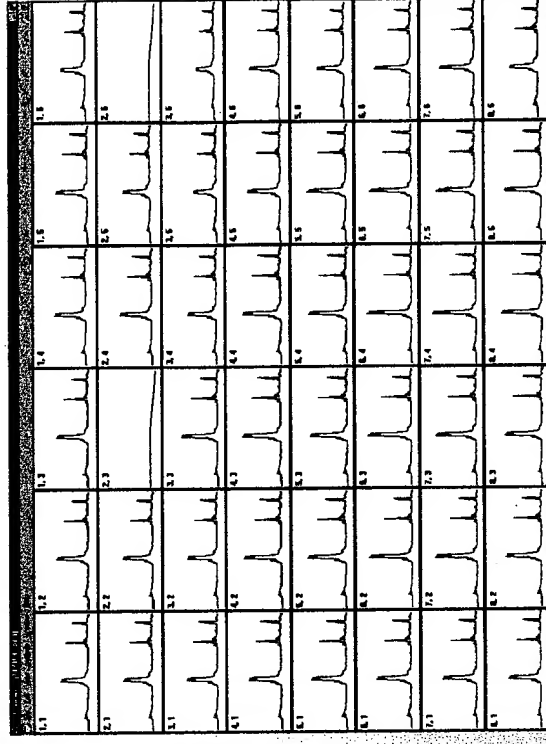
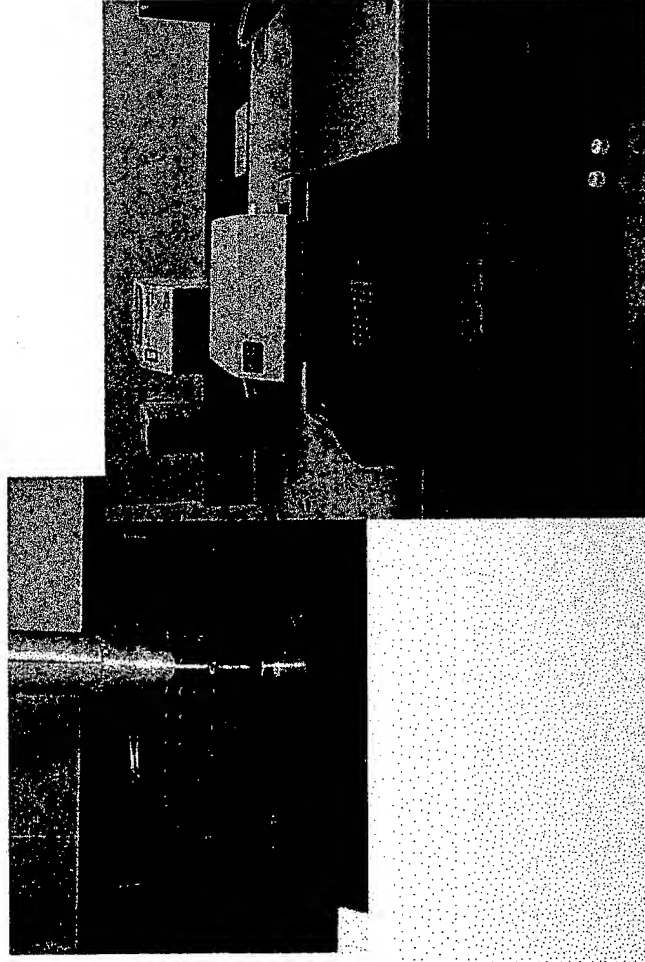
- 48 reactions per run
- 6 mL reaction volume
- Reagents injected robotically at T and P
- Real-time monitoring
 - T, P and gas uptake in cell

Secondary Screening Analysis: FT-IR



IR spectra of polymers allows rapid analysis of:

- co-monomer composition
- tacticity

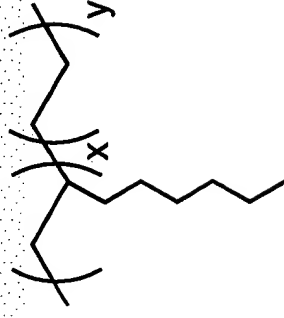


Automated Reflection FT-IR Analysis of Polymer Films

A Case Study of High-Throughput Catalyst Discovery

Target: High Temperature Solution Phase LLDPE Catalyst

- ❖ Catalyst must operate at temperatures > 130 °C
- ❖ Requires high activity, high M_w , high 1-octene incorporation



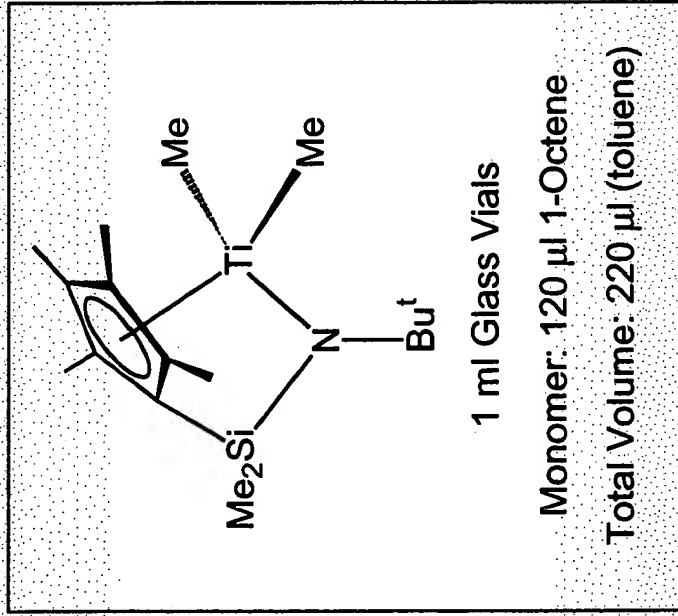
Primary Screen:
1-Octene Polymerization Activity

Hits and Information

1-Octene Polymerization Capability

Secondary Screen:
High Temperature Ethylene/ 1-Octene Screen

Adaption to Small Scale: Primary Screening Validation

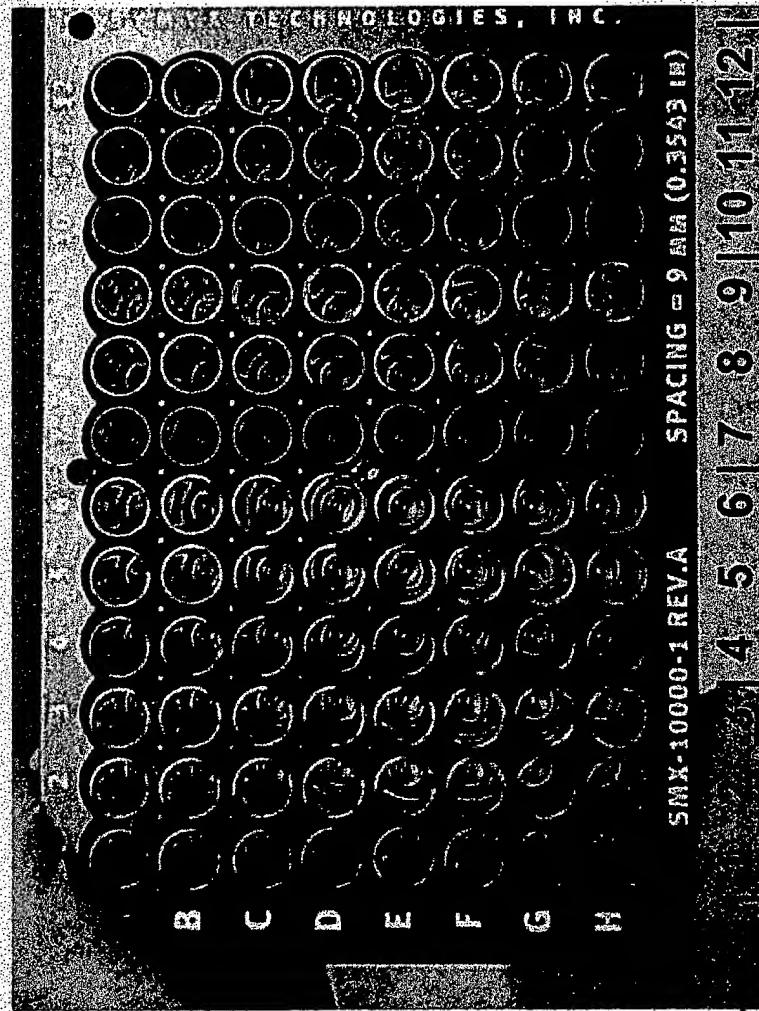
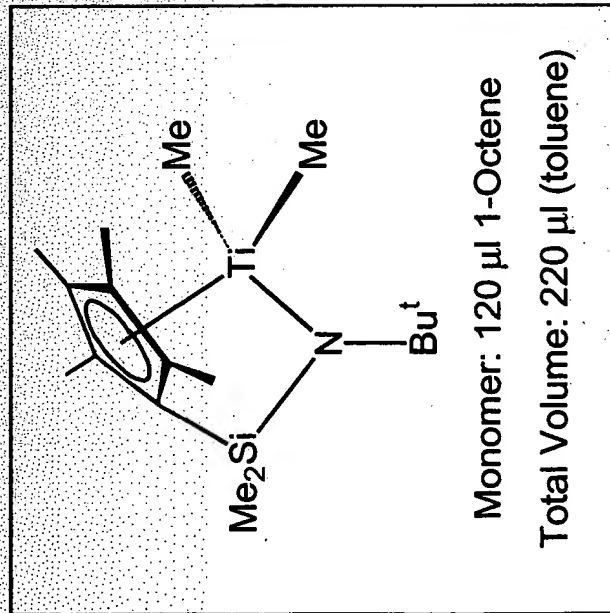


[8 x 12] Array

3 Catalyst Concentrations
4 Activation Methods
8 Replicates of Each
Condition

Use robots to dispense
reagents and to analyze
products for speed, accuracy

Adaption to Small Scale: Primary Screening Validation



8
Replicates

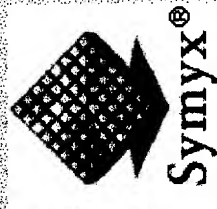
Column No.

4 Activation
Zones

3 Catalyst
Concentrations
(mM)

| None (Control) | | | | 200 eq. MAO | | | | [Me ₂ PhNH] ⁺ [B(C ₆ F ₅) ₄] ⁻ | | | | [Me ₂ PhNH] ⁺ [B(C ₆ F ₅) ₄] ⁻ + 5 eq. Bu ⁱ ₃ Al | | | |
|-------------------|------|-------|-----|----------------|-------|-------|-----|---|------|-------|-----|---|-------|-----|-------|
| 1.8 | 0.18 | 0.018 | 1.8 | 0.18 | 0.018 | 0.018 | 1.8 | 1.8 | 0.18 | 0.018 | 1.8 | 0.18 | 0.018 | 1.8 | 0.018 |

Adaption to Small Scale: Primary Screening Validation Procedure

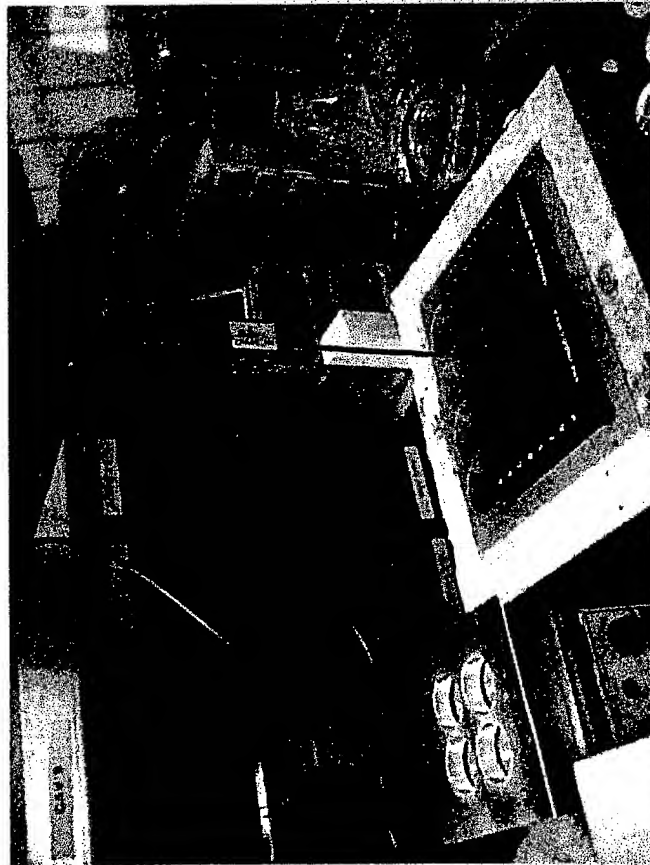


1. Add 120 μL 1-Octene.
2. Add Bu^i_3Al to columns 10, 11, 12.
3. Add $[(\eta^5\text{-C}_5\text{Me}_5)\text{SiMe}_2(\eta^1\text{-NBu}^i)\text{TiMe}_2]$.
4. Add Activators.



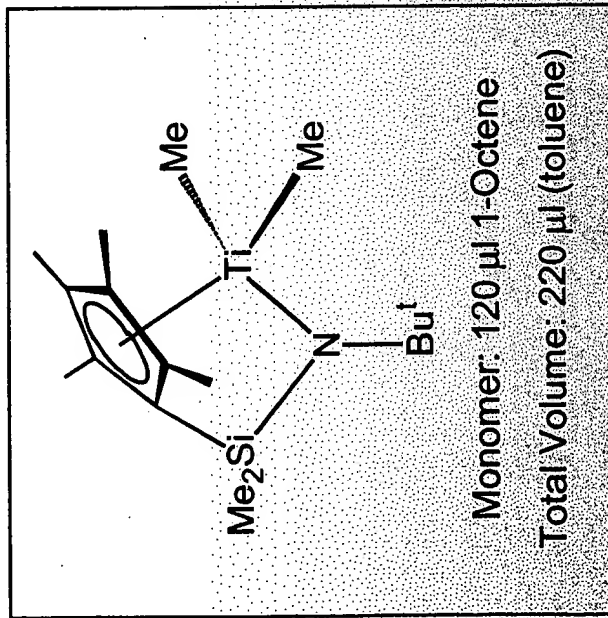
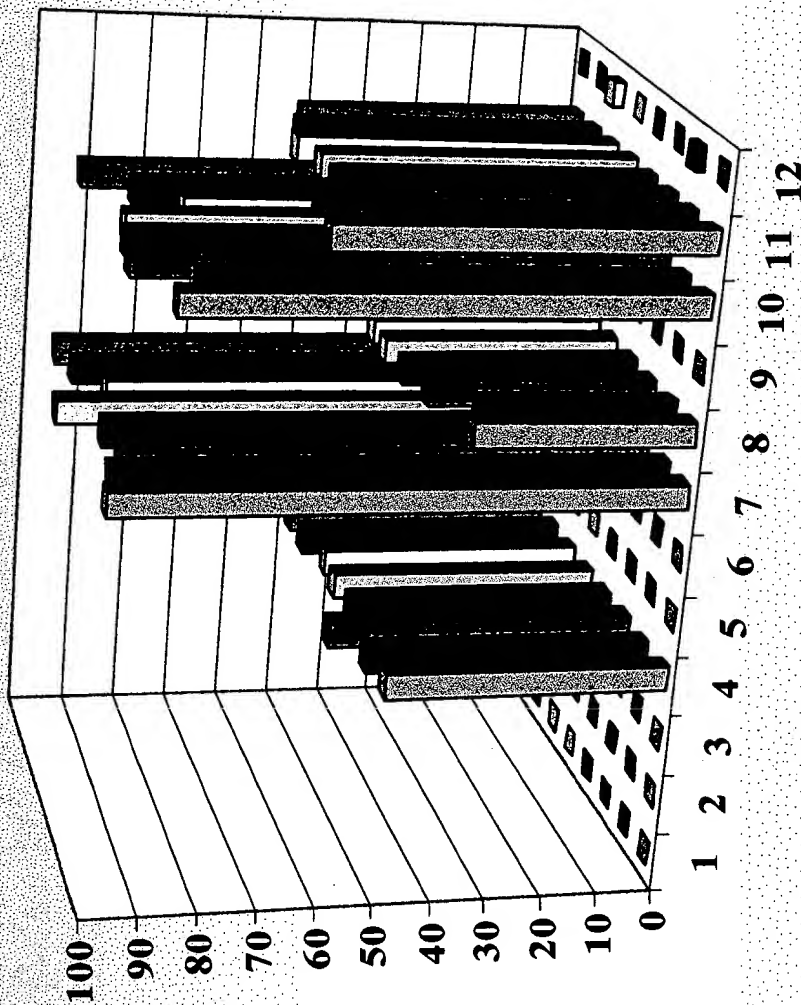
1. Run Polymerization
2. Quench, Filter
3. Automated Dilution
4. Rapid GPCTM
5. Obtain M_w , Conversion

All Procedures Employ
Robotic Liquid Handling



Adaption to Small Scale: Primary Screening Catalyst Activity Validation

% Conversion
to Poly-1-Octene

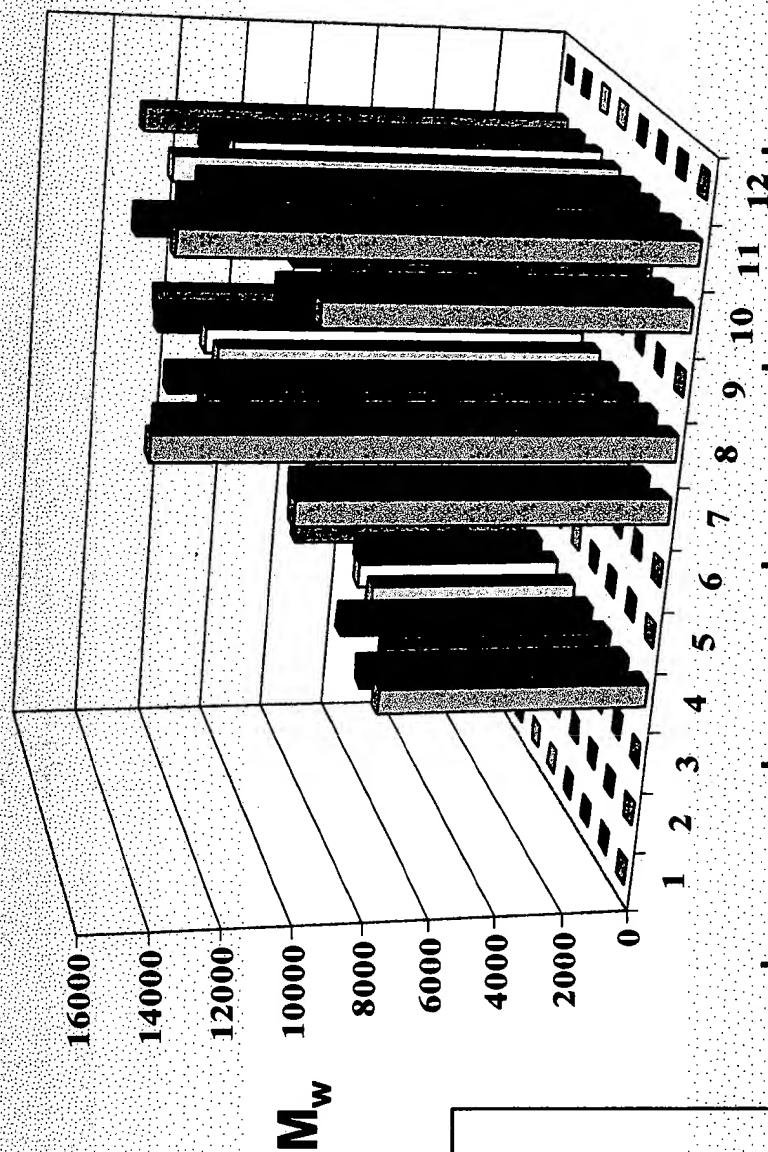


| None | 200 eq. MAO | [Me ₂ PhNH] ⁺ [B(C ₆ F ₅) ₄] ⁻ | | | [Me ₂ PhNH] ⁺ [B(C ₆ F ₅) ₄] ⁻ + 5 eq. Bu ^t Al | | |
|------|----------------|---|-----|------|---|------|------|
| 1.8 | 0.18 | 0.18 | 1.8 | 0.18 | 1.8 | 0.18 | 0.18 |

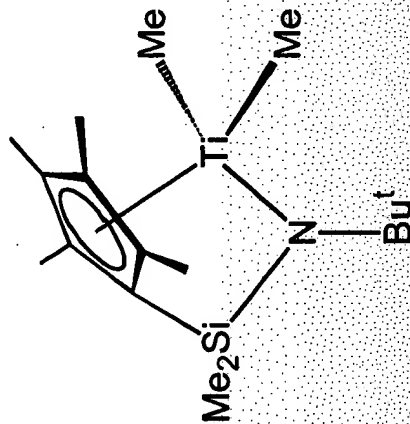
3 Catalyst
Concentrations
(mM)

Adaption to Small Scale: Primary Screening

Mw Validation



M_w



Monomer: 120 μl 1-Octene
Total Volume: 220 μl (toluene)

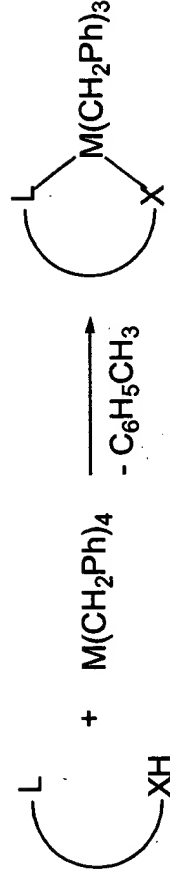
| None | | | 200 eq. MAO | | | [Me ₂ PhNH] ⁺ [B(C ₆ F ₅) ₄] ⁻ | | | [Me ₂ PhNH] ⁺ [B(C ₆ F ₅) ₄] ⁻ + 5 eq. Bu ^t Al | | |
|------|------|------|-------------|------|------|---|------|------|---|------|------|
| 1.8 | 0.18 | 0.18 | 1.8 | 0.18 | 0.18 | 1.8 | 0.18 | 0.18 | 1.8 | 0.18 | 0.18 |
| | | | | | | | | | | | |

3 Catalyst
Concentrations
(mM)

Catalyst Discovery Screening

- Adaption of the Chemistry to Smaller Scale
- Large Number of Ligands with Suitable Diversity
- Efficient Methods of Attaching Ligands to Metals
- Suitable Rapid Screening Technique

The Use of Acidic ligands with $M(\text{CH}_2\text{Ph})_4$ ($M = \text{Zr, Hf}$) and Multiple Activation Conditions



$\text{M}(\text{CH}_2\text{Ph})_4$
($M = \text{Zr, Hf}$)

[3 x 8] Set of
Charged Ligands

1 Hour, Toluene

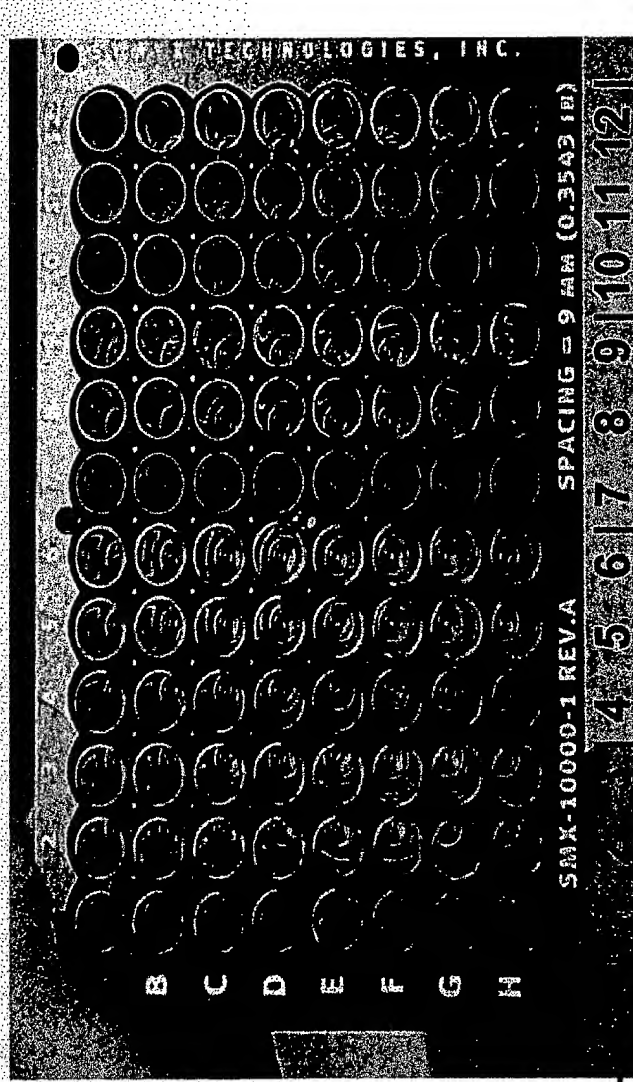
Ligand/ $\text{M}(\text{CH}_2\text{Ph})_4$
Mixtures

Multiple Activation
Conditions

Catalyst Discovery Screening - Plate Design

Discovery Plate Design

Eight Activation
Conditions for Each
Ligand/ $M(CH_2Ph)_4$
Combination
- 2 plates of 4
conditions



Column No.

Activation
Zones

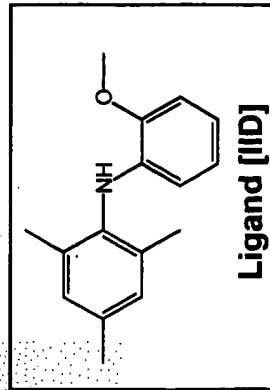
Ligand/ $M(CH_2Ph)_4$
Concentration
(mM)

| 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 | Row No. | | | |
|--|---|---|---|--|---|---|---|---|----|----|----|---|--|--|--|
| $[Ph_3C]^+$ $[B(C_6F_5)_4]^-$ 5 eq. Et_3Al | | | | $[Ph_3C]^+$ $[B(C_6F_5)_4]^-$ 5 eq. Bu^i_3Al | | | | $[Me_2PhNH]^+$ $[B(C_6F_5)_4]^-$ 5 eq. Et_3Al | | | | $[Me_2PhNH]^+$ $[B(C_6F_5)_4]^-$ 5 eq. Bu^i_3Al | | | |
| 1.6 | | | | 1.6 | | | | 1.6 | | | | 1.6 | | | |

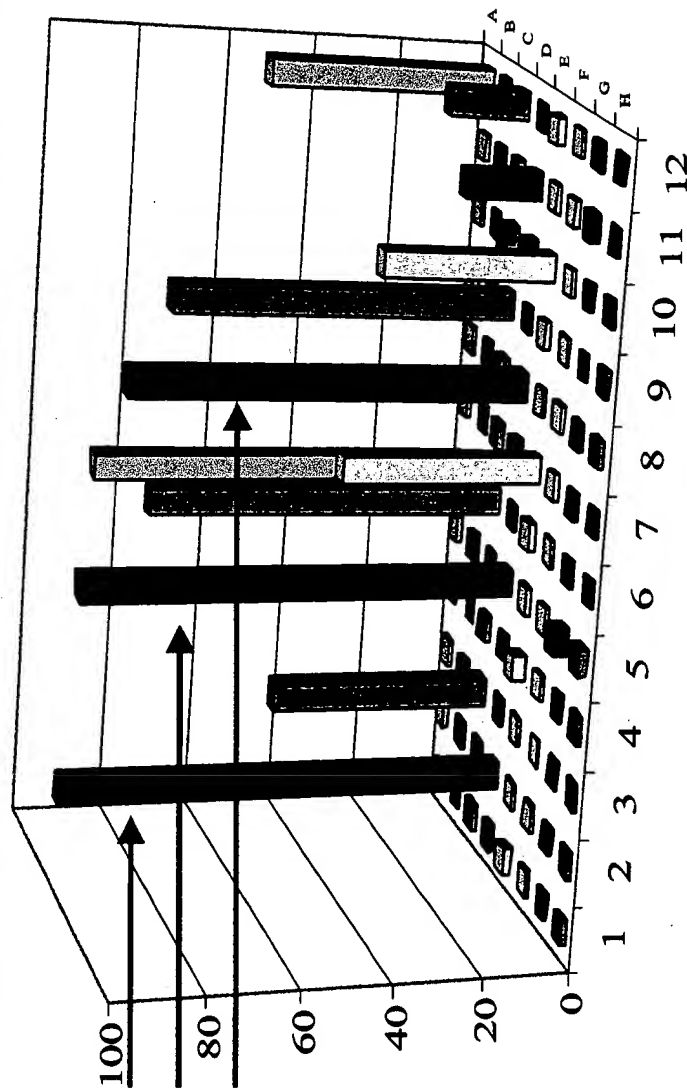
4 Zones each containing [3 x 8] Ligand Set

Catalyst Discovery Screening - Activity Results

% Conversion Data
Hf(CH₂Ph)₄



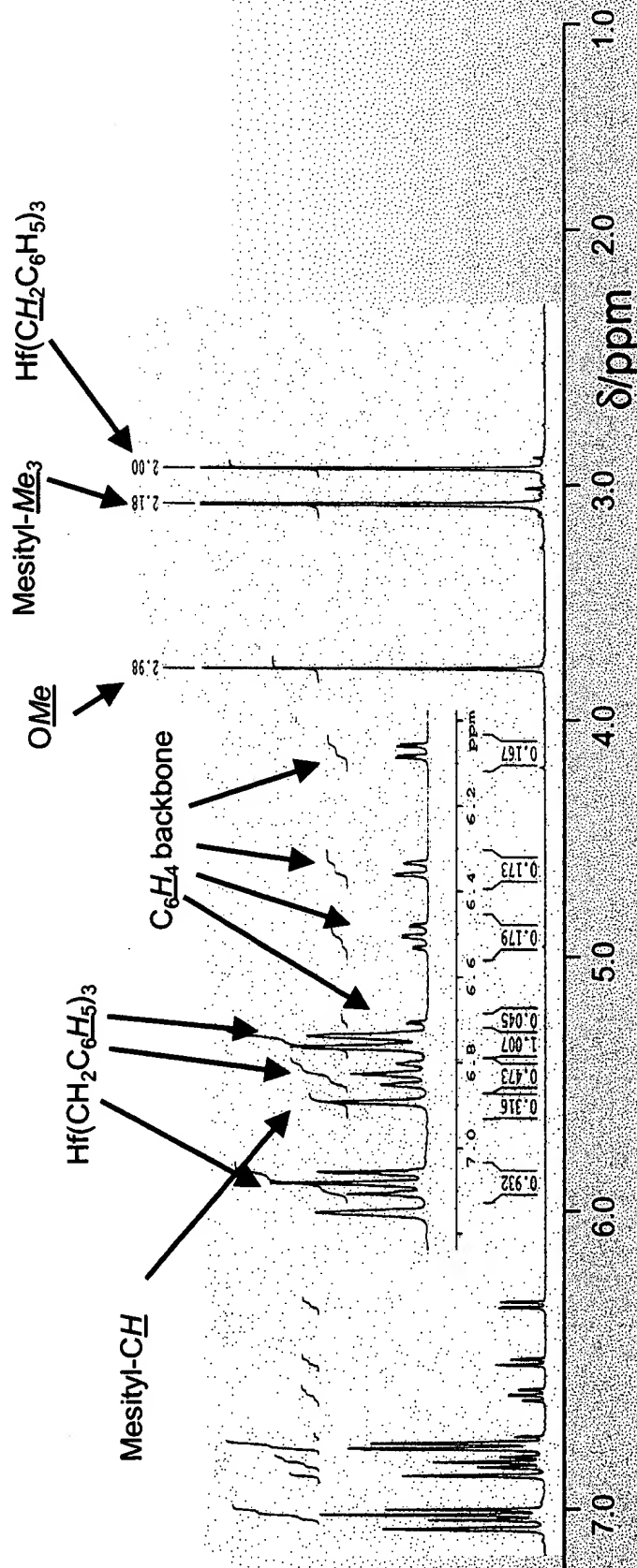
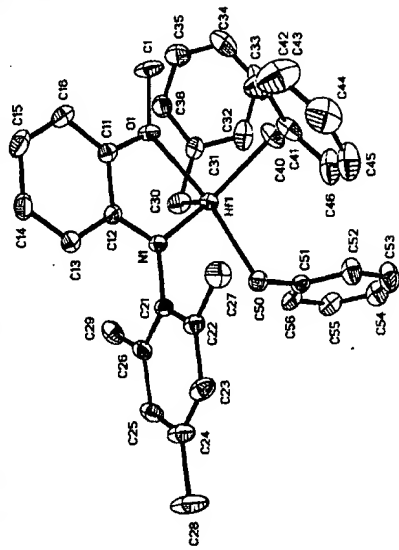
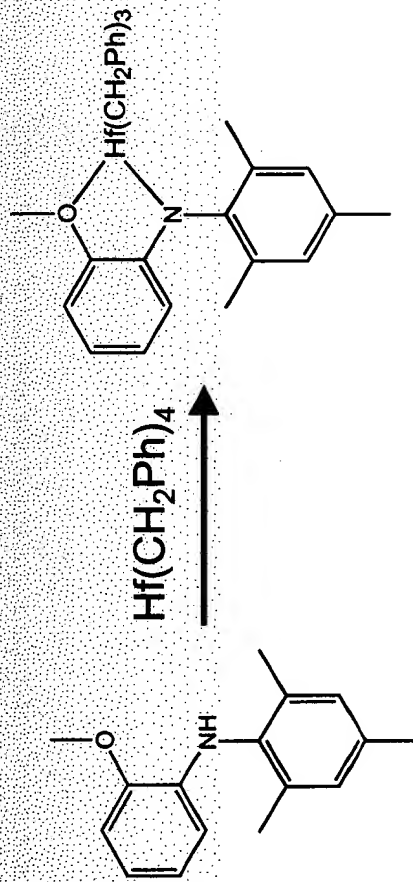
% Conversion
to Poly-1-Octene



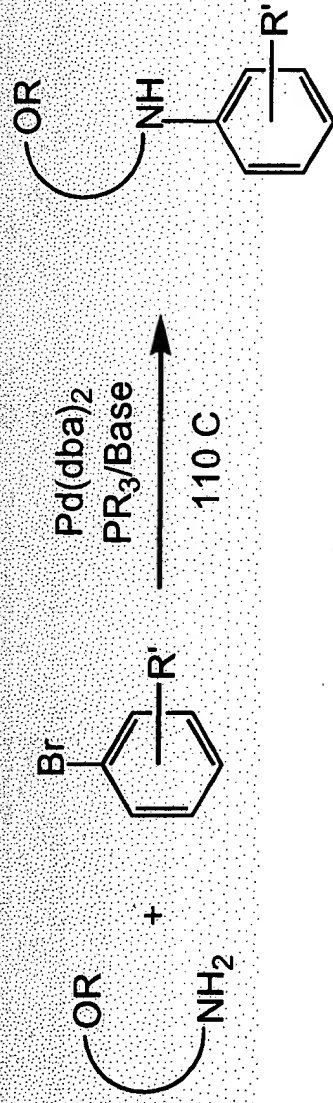
Ligand/Hf(CH₂Ph)₄
Concentration
(mM)

| | | | |
|--|---|---|--|
| [Me ₂ PhNH] ⁺ [B(C ₆ F ₅) ₄] ⁻ + 5 eq. Et ₃ Al | [Me ₂ PhNH] ⁺ [B(C ₆ F ₅) ₄] ⁻ + 5 eq. Bu ₁₃ Al | [Ph ₃ C] ⁺ [B(C ₆ F ₅) ₄] ⁻ + 5 eq. Et ₃ Al | [Ph ₃ C] ⁺ [B(C ₆ F ₅) ₄] ⁻ + 5 eq. Bu ¹ ₃ Al |
| 1.6 | 1.6 | 1.6 | 1.6 |

After Hit Identification, Conventional Chemistry Identifies Structure of Active Catalyst

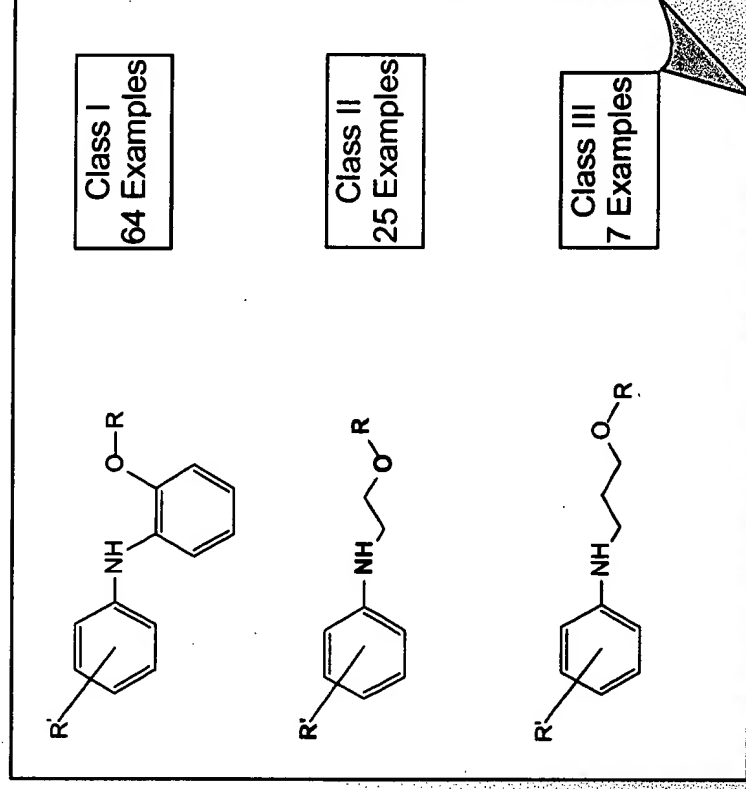
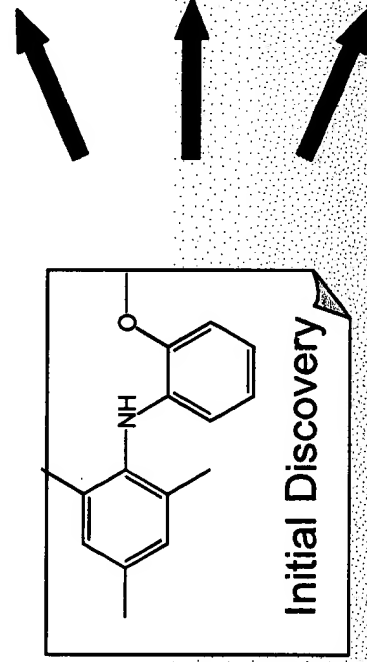


Focus Library of Structural Variation Around Identified Hit

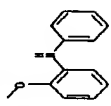
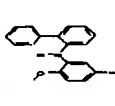
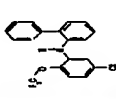
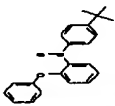
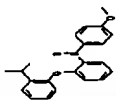
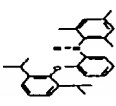
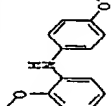
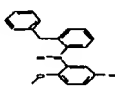
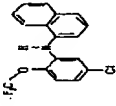
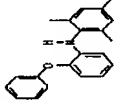
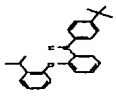
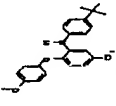
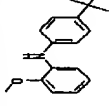
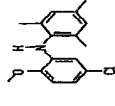
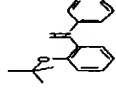
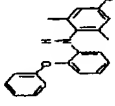
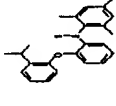
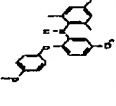
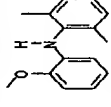
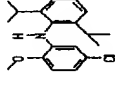
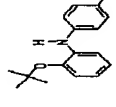
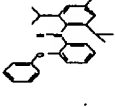
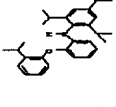
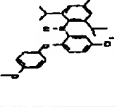
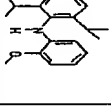
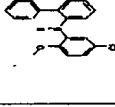
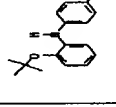
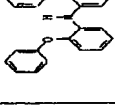
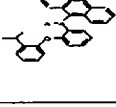

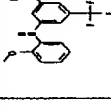
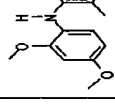
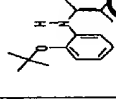
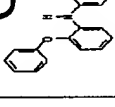
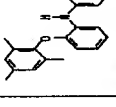
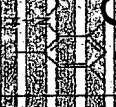
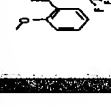
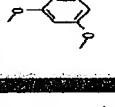
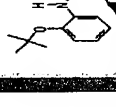
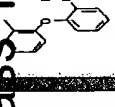
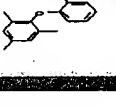

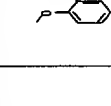
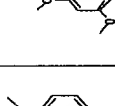
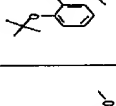
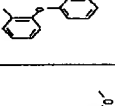
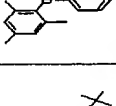

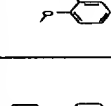
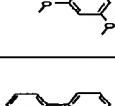
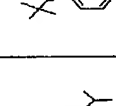
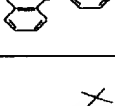
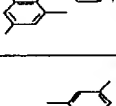

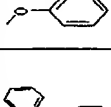
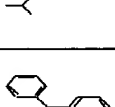
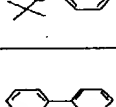
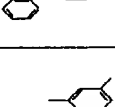
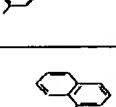
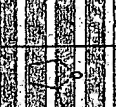
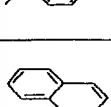
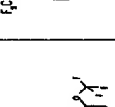
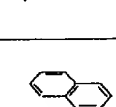
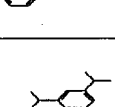
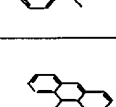

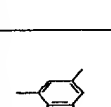
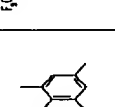
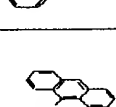
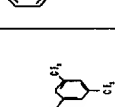
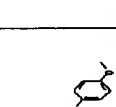
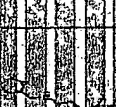
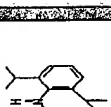
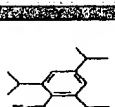
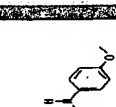
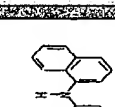
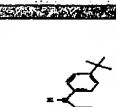

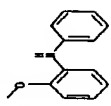
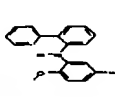
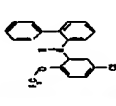
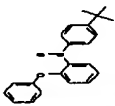
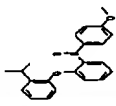
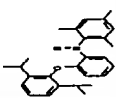
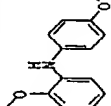
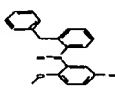
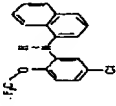
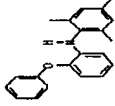
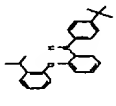
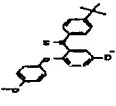
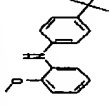
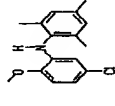
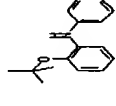
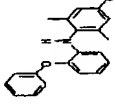
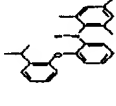
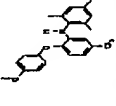
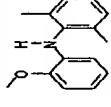
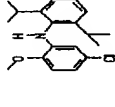
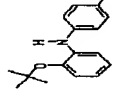
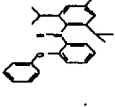
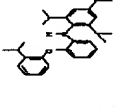
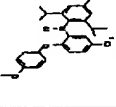
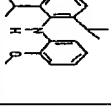
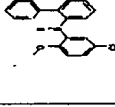
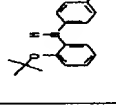
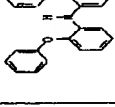
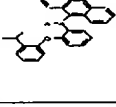

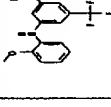
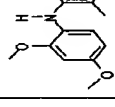
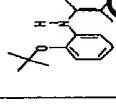
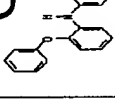
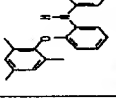
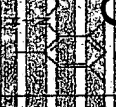
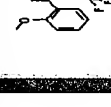
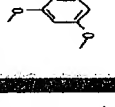
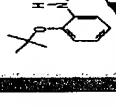
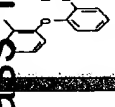
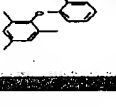

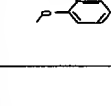
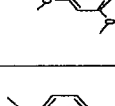
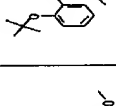
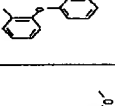
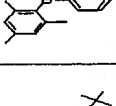

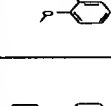
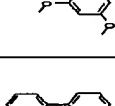
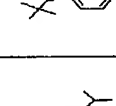
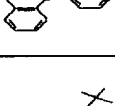
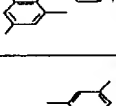

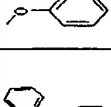
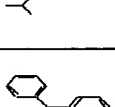
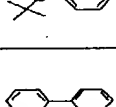
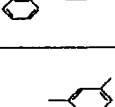
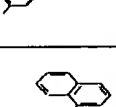
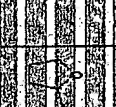
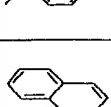
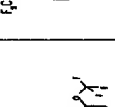
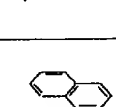
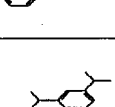
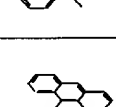

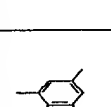
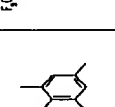
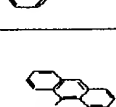
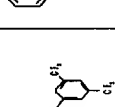
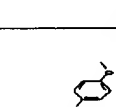
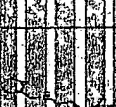
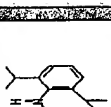
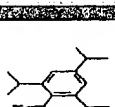
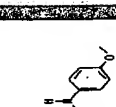
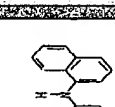
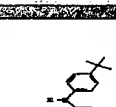

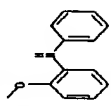
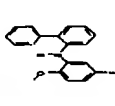
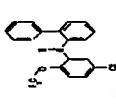
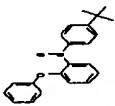
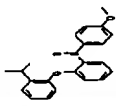
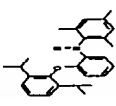
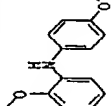
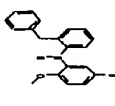
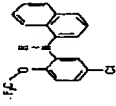
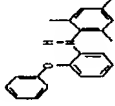
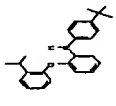
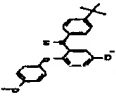
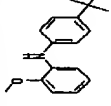
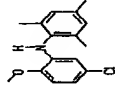
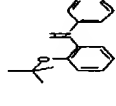
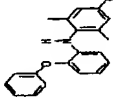
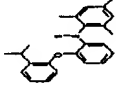
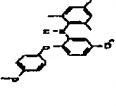
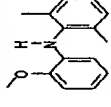
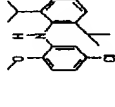
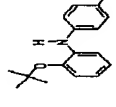
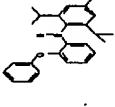
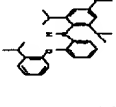
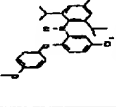
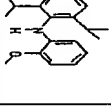
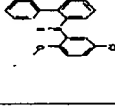
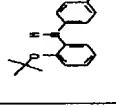
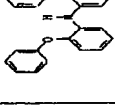
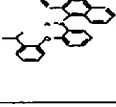

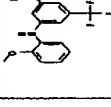
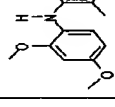
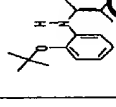
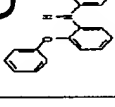
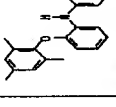
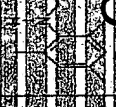
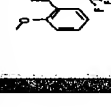
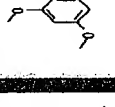
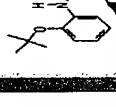
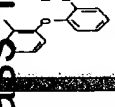
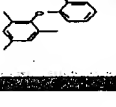

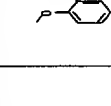
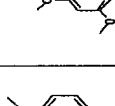
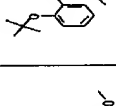
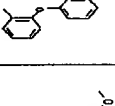
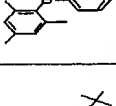

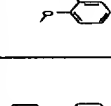
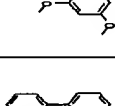
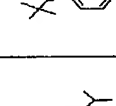
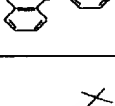
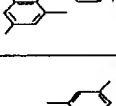

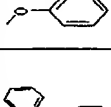
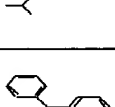
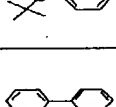
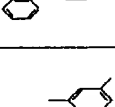
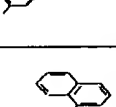
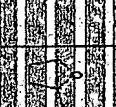
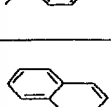
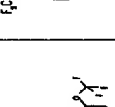
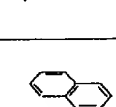
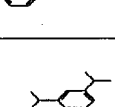
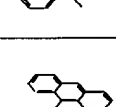

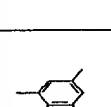
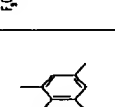
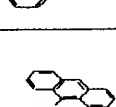
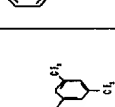
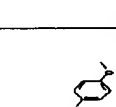
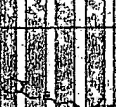
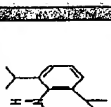
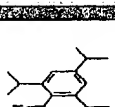
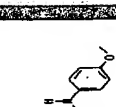
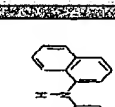
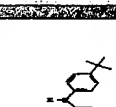



Focus Library Preparation

❖ 96 Structural Variations



Focus Library for Amine Ether Hit

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| Class 3 | | | | | |

Secondary Screening

96 High Temperature Ethylene-Octene Copolymerizations

Reactor Conditions

130 °C, 100 psi Ethylene

6.1 mL Total (Toluene)

0.25 mL 1-Octene

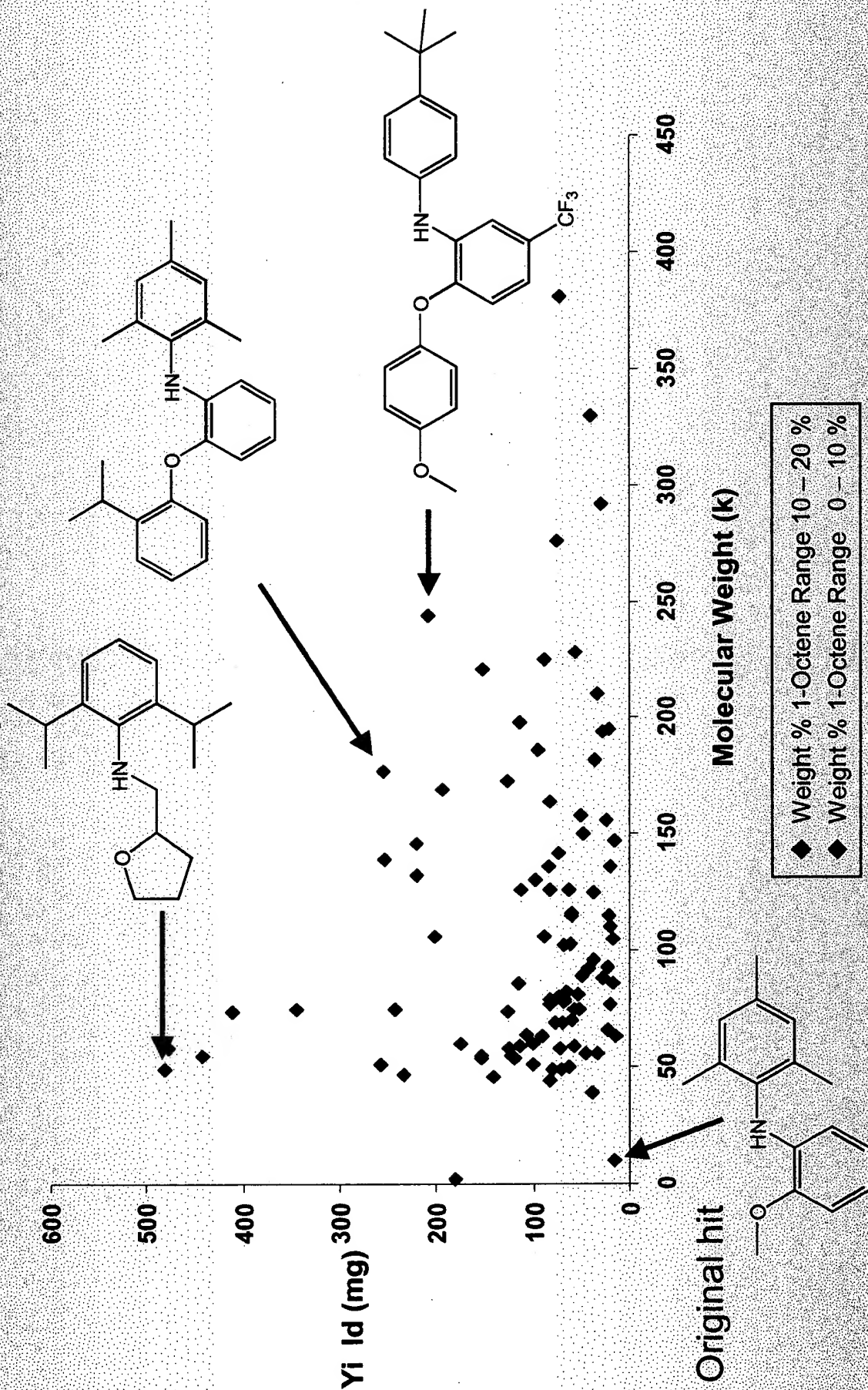
2 μmol Bu^i_3Al

1 μmol $\text{Hf}(\text{CH}_2\text{C}_6\text{H}_5)_4/\text{Ligand}$

Target:
200-300 mg Copolymer

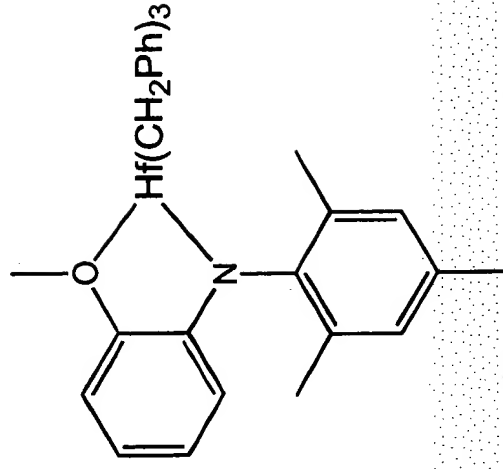
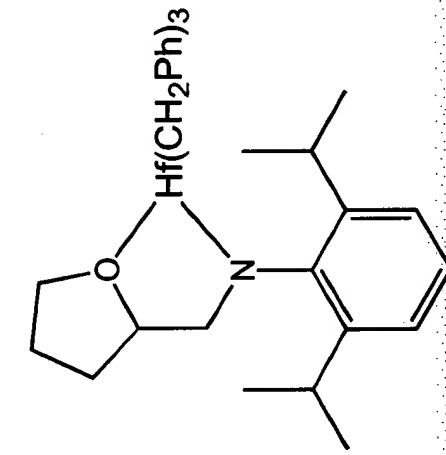
- | | | | |
|---------------------|---|---|----------------|
| [8 x 12] Ligand Set | $\xrightarrow{\text{Hf}(\text{CH}_2\text{Ph})_4}$ | 1) Add Bu^i_3Al | Obtain yield, |
| | | 2) Add $[\text{Me}_2\text{PhNH}][\text{B}(\text{C}_6\text{F}_5)_4]$ | molecular |
| | $\xrightarrow{1 \text{ Hour, Toluene}}$ | 3) Inject Ligand/ $\text{Hf}(\text{CH}_2\text{Ph})_4$ | weight, MWD, |
| | | 4) Monitor Run for 30 Minutes | wt. % 1-octene |

Focus Library Identifies Large Structure / Activity Effects in this Family, Improved Catalysts



1 Gallon Industrial Ethylene 1-Octene Autoclave Results Using Isolated Catalysts Validate Screening Results

Performed in Solution
Phase
at 130 °C



| | | |
|---------------------------|--------|-------|
| Activity (g poly/mmol Hf) | 13,760 | 7,560 |
| M_w | 181 k | 97 k |
| M_w/M_n | 2.8 | 3.0 |
| Density | 0.917 | 0.910 |

Conclusions

Effective high-throughput combinatorial methods have been developed for organometallic polyolefin catalyst systems

Months or years of conventional catalyst research can be accomplished in days or weeks

Small-scale 1-octene polymerization screens were validated using a commercial catalyst

New hafnium post-metallocene catalyst family was discovered

Focus library of 96 structural variations was rapidly investigated

- significant enhancements observed
- wide variation in structure / property response

Optimized structures scaled up to commercial-scale process reactors